## Extending Ismeans

Russell V. Lenth

March 14, 2015

## 1 Introduction

Suppose you want to use Ismeans for some type of model that it doesn't (yet) support. Or, suppose you have developed a new package with a fancy model-fitting function, and you'd like it to work with Ismeans. What can you do? Well, there is hope because Ismeans is designed to be extended.

The first thing to do is to look at the help page for extending the package:

```
R> help("extending-lsmeans", package="lsmeans")
```

It gives details about the fact that you need to write two S3 methods, recover.data and lsm.basis, for the class of object that your model-fitting function returns. The recover.data method is needed to recreate the dataset so that the reference grid can be identified. The lsm.basis method then determines the linear functions needed to evaluate each point in the reference grid and to obtain associated information—such as the variance-covariance matrix—needed to do estimation and testing.

This vignette presents an example where suitable methods are developed, and discusses a few issues that arise.

# 2 Data example

The MASS package contains various functions that do robust or outlier-resistant model fitting. We will cobble together some lsmeans support for these. But first, let's create a suitable dataset (a simulated two-factor experiment) for testing.<sup>1</sup>

```
R> fake = expand.grid(rep = 1:5, A = c("a1", "a2"), B = c("b1", "b2", "b3"))
R> fake$y = c(11.46,12.93,11.87,11.01,11.92,17.80,13.41,13.96,14.27,15.82,
23.14,23.75,-2.09,28.43,23.01,24.11,25.51,24.11,23.95,30.37,
17.75,18.28,17.82,18.52,16.33,20.58,20.55,20.77,21.21,20.10)
```

The y values were generated using predetermined means and Cauchy-distributed errors. There are some serious outliers in these data.

## 3 Supporting rlm

The MASS package provides an rlm function that fits robust-regression models using M estimation. We'll fit a model using the default settings for all tuning parameters:

 $<sup>^{1}\</sup>mathrm{I}$  unapologetically use = as the assignment operator. It is good enough for C and Java, and supported by R.

```
R> library(MASS)
R > fake.rlm = rlm(y ~A * B, data = fake)
R> library(lsmeans)
R> lsmeans(fake.rlm, ~B | A)
A = a1:
 В
     lsmean
                    SE df asymp.LCL asymp.UCL
b1 11.83800 0.4774474 NA
                          10.90222 12.77378
b2 23.30000 0.4774474 NA
                          22.36422
                                     24.23578
b3 17.80078 0.4774474 NA 16.86500
                                    18.73656
A = a2:
 В
     lsmean
                    SE df asymp.LCL asymp.UCL
b1 14.68344 0.4774474 NA
                          13.74766
                                    15.61922
b2 24.71164 0.4774474 NA 23.77586
                                    25.64742
b3 20.64200 0.4774474 NA 19.70622 21.57778
Confidence level used: 0.95
```

The first lesson to learn about extending Ismeans is that sometimes, it already works! It works here because rlm objects inherit from lm, which is supported by the Ismeans package, and rlm objects aren't enough different to create any problems.

## 4 Supporting lqs objects

The MASS resistant-regression functions lqs, lmsreg, and ltsreg are another story, however. They create lqs objects that are not extensions of any other class, and have other issues, including not even having a vcov method. So for these, we really do need to write new methods for lqs objects. First, let's fit a model.

```
R> fake.lts = ltsreg(y ~ A * B, data = fake)
```

#### 4.1 The recover data method

It is usually an easy matter to write a recover.data method. Look at the one for lm objects:

Note that all it does is obtain the call component and call the method for class "call", with additional arguments for its terms component and na.action. It happens that we can access these attributes in exactly the same way as for lm objects; so, ...

```
R> recover.data.lqs = lsmeans:::recover.data.lm
```

The trickier part is testing it, as it isn't clear that there is a required data argument.

```
R> rec.fake = recover.data(fake.lts, data = NULL)
R> head(rec.fake)

A B
1 a1 b1
2 a1 b1
3 a1 b1
4 a1 b1
5 a1 b1
6 a2 b1
```

Our recovered data excludes the response variable y (owing to the delete.response call), and this is fine.

By the way, the data argument is handed to recover.data if it is specified in the ref.grid or lsmeans call. It is needed to cover a desperate situation that occurs with certain kinds of models that are fitted by iteratively modifying the data. In those cases, the only way to recover the data is to for the user to give it explicitly, and recover.data just adds a few needed attributes to it.

Related (in a surprising way) to the above, if you check for any error conditions in recover.data, simply have it return a character string with the desired message, rather than invoking stop. This provides a cleaner exit. The reason is that whenever recover.data throws an error, an informative message suggesting that data be provided is displayed. But a character return value is tested for an throws a different error with your message.

#### 4.2 The lsm.basis method

The lsm.basis method has four required arguments:

```
R> args(lsmeans:::lsm.basis.lm)
function (object, trms, xlev, grid, ...)
NULL
```

These are, respectively, the model object, its terms component (at least for the right-hand side of the model), a list of levels of the factors, and the grid of predictor combinations that specify the reference grid.

The function must obtain six things and return them in a named list. They are the matrix X of linear functions for each point in the reference grid, the regression coefficients bhat; the variance-covariance matrix V; a matrix nbasis for non-estimable functions; a function dffun(k,dfargs) for computing degrees of freedom for the linear function sum(k\*bhat); and a list dfargs of arguments to pass to dffun.

To write your own lsm.basis function, examining some of the existing methods can help; but the best resource is the predict method for the object in question, looking carefully to see what it does to predict values for a new set of predictors (e.g., newdata in predict.lm). Following this advice, let's take a look at it:

```
R> MASS:::predict.lqs
```

```
function (object, newdata, na.action = na.pass, ...)
   {
       if (missing(newdata))
           return(fitted(object))
       Terms <- delete.response(terms(object))</pre>
       m <- model.frame(Terms, newdata, na.action = na.action, xlev = object$xlevels)</pre>
       if (!is.null(cl <- attr(Terms, "dataClasses")))</pre>
            .checkMFClasses(cl, m)
       X <- model.matrix(Terms, m, contrasts = object$contrasts)</pre>
       drop(X %*% object$coefficients)
   <bytecode: 0x00000000bf0b3a0>
   <environment: namespace:MASS>
   Based on this, here is a listing of an lsm.basis method for lqs objects:
   R> lsm.basis.lqs = function(object, trms, xlev, grid, ...) {
          m = model.frame(trms, grid, na.action = na.pass, xlev = xlev)
2
          X = model.matrix(trms, m, contrasts.arg = object$contrasts)
3
          bhat = coef(object)
          Xmat = model.matrix(trms, data=object$model)
          V = rev(object$scale)[1]^2 * solve(t(Xmat) %*% Xmat)
          nbasis = matrix(NA)
          dfargs = list(df = nrow(Xmat) - ncol(Xmat))
8
          dffun = function(k, dfargs) dfargs$df
          list(X=X, bhat=bhat, nbasis=nbasis, V=V, dffun=dffun, dfargs=dfargs)
10
      }
11
   Before explaining it, let's verify that it works:
   R> lsmeans(fake.lts, ~ B | A)
   A = a1:
                        SE df lower.CL upper.CL
    b1 11.87278 0.2284451 24 11.40129 12.34427
    b2 23.09278 0.2284451 24 22.62129 23.56427
    b3 17.77278 0.2284451 24 17.30129 18.24427
   A = a2:
         lsmean
                        SE df lower.CL upper.CL
    b1 13.91278 0.2284451 24 13.44129 14.38427
    b2 24.06278 0.2284451 24 23.59129 24.53427
    b3 20.50278 0.2284451 24 20.03129 20.97427
   Confidence level used: 0.95
```

Hooray! Note the results are comparable to those we had for fake.rlm, albeit the standard errors are quite a bit smaller.

### 4.3 Dissecting lsm.basis.lqs

Let's go through the listing of this method, by line numbers.

- 2-3: Construct the linear functions, X. This is a pretty standard standard two-step process: First obtain a model frame, m, for the grid of predictors, then pass it as data to model.data to create the associated design matrix. As promised, this code is essentially identical to what you find in predict.lqs.
  - 4: Obtain the coefficients, bhat. Most model objects have a coef method.
- 5–6: Obtain the covariance matrix, V, of bhat. In many models, this can be obtained using the object's vcov method. But not in this case. Instead, I cobbled one together using what it would be for ordinary regression:  $\hat{\sigma}^2(X'X)^{-1}$ , where X is the design matrix for the whole dataset (not the reference grid). Here,  $\hat{\sigma}$  is obtained using the last element of the scale element of the object (depending on the method, there are one or two scale estimates). This probably under-estimates the variances and distorts the covariances, because robust estimators have some efficiency loss.
  - 7: Compute the basis for non-estimable functions. This applies only when there is a possibility of rank deficiency in the model, and lqs methods cannot handle that. All linear functions are estimable, and we signal that by setting nbasis equal to a 1 × 1 matrix of NA. If rank deficiency were possible, the estimability package (which is required by Ismeans) provides a nonest.basis function that makes this fairly painless—I would have coded:

```
R> nbasis = estimability::nonest.basis(Xmat)
```

On the other hand, if rank-deficient cases are not possible, set nbasis equal to all.estble, a constant in the estimability package.

There is a subtlety you need to know regarding estimability. Suppose the model is rank-deficient, so that the design matrix  $\mathbf{X}$  has p columns but rank r < p. In that case, bhat should be of length p (not r), and there should be p-r elements equal to NA, corresponding to columns of  $\mathbf{X}$  that were excluded from the fit. Also,  $\mathbf{X}$  should have all p columns. In other words, do not alter or throw-out columns of  $\mathbf{X}$  or their corresponding elements of bhat—even those with NA coefficients—as they are essential for assessing estimability.  $\mathbf{V}$  should be  $r \times r$ , however: the covariance matrix for the non-excluded predictors.

- 8-9: Obtain dffun and dfargs. This is a little awkward because it is designed to allow support for mixed models, where approximate methods may be used to obtain degrees of freedom. The function dffun is expected to have two arguments: k, the vector of coefficients of bhat, and dfargs, a list containing any additional arguments. In this case (and in many other models), the degrees of freedom are the same regardless of k. We put the required degrees of freedom in dfargs and write dffun so that it simply returns that value.
- 10: Return these results in a named list.

#### 4.4 The "honest" version

Because of the inadequacies mentioned above for estimating the covariance matrix, then—lacking any better estimate—I think it's probably better to set it and the degrees of freedom to NAs. We will still be able to get the LS means and contrasts thereof, but no standard errors or tests. With that in mind, here's a replacement version:

```
R> lsm.basis.lqs = function(object, trms, xlev, grid, ...) {
       m = model.frame(trms, grid, na.action = na.pass, xlev = xlev)
       X = model.matrix(trms, m, contrasts.arg = object$contrasts)
       bhat = coef(object)
       V = diag(rep(NA, length(bhat)))
       nbasis = matrix(NA)
       dffun = function(k, dfargs) NA
       list(X=X, bhat=bhat, nbasis=nbasis, V=V, dffun=dffun, dfargs=list())
  }
And here is a test:
R> lsmeans(fake.lts, pairwise ~ B)
$1smeans
      1smean SE df asymp.LCL asymp.UCL
b1 12.89278 NA NA
                          NA
b2 23.57778 NA NA
                           NΑ
                                     NA
b3 19.13778 NA NA
                          NA
                                     NA
Results are averaged over the levels of: A
Confidence level used: 0.95
$contrasts
 contrast estimate SE df z.ratio p.value
b1 - b2
           -10.685 NA NA
                               NA
                                       NA
b1 - b3
            -6.245 NA NA
                               NA
                                       NA
b2 - b3
             4.440 NA NA
                               NA
                                       NA
Results are averaged over the levels of: A
P value adjustment: tukey method for comparing a family of 3 estimates
```

### 5 Hook functions

Most linear models supported by Ismeans have straightforward structure: Regression coefficients, their covaraince matrix, and a set of linear functions that define the reference grid. However, a few are more complex. An example is the "clm" class in the ordinal package, which allows a scale model in addition to the location model. When a scale model is used, the scale parameters are included in the model matrix, regression coefficients, and covariance matrix, and we can't just use the usual matrix operations to obtain estimates and standard errors. To facilitate using custom routines for these tasks, the lsm.basis.clm function includes, in its misc part, the names (as character constants) of two "hook" functions: misc\$estHook has the name of the function to call when computing estimates, standard errors, and degrees of freedom (for thesummary method); and misc\$vcovHook has the name of the function to call to obtain the covariance matrix of the grid values (used by the vcov method). These functions are called in lieu of the usual built-in routines for these purposes, and return the appropriately sized matrices.

In addition, you may want to apply some form of special post-processing after the reference grid is constructed. To provide for this, give the name of your function to post-process the object

in misc\$postGridHook. Again, "clm" objects (as well as "polr" in the MASS package) serve as an example. They allow a mode specification that in twio cases, calls for post-processing. The "cum.prob" mode uses the regrid function to transform the linear predictor to the cumulative-probability scale. And the "prob" mode performs this, as well as applying the contrasts necessary to difference the cumulative probabilities into the class probabilities.

## 6 Conclusions

It is relatively simple to write appropriate methods that work with Ismeans for model objects it does not support. I hope this vignette is helpful for understanding how. Furthermore, if you are the developer of a package that fits linear models, I encourage you to include recover.data and lsm.basis methods for those classes of objects, and to remember to export them in your NAMESPACE file as follows:

S3method(myobject, recover.data)
S3method(myobject, lsm.basis)